

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
 Data File : BP013530.D
 Acq On : 18 Jan 2023 19:55
 Operator : CG/JU
 Sample : 01146-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43Z9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013530.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.293	15	18	29	rVB	102482	146136	1.70%	0.151%
2	3.758	95	97	104	rBV2	26548	46608	0.54%	0.048%
3	3.822	105	108	117	rVB	94585	133131	1.55%	0.137%
4	3.946	125	129	135	rVB	86138	133484	1.56%	0.138%
5	4.046	142	146	152	rVB	36432	54584	0.64%	0.056%
6	4.152	160	164	169	rVB	58676	79763	0.93%	0.082%
7	4.275	181	185	191	rVB3	35729	50942	0.59%	0.053%
8	4.422	207	210	215	rVB7	20953	31570	0.37%	0.033%
9	4.611	239	242	247	rVB3	32541	40275	0.47%	0.042%
10	4.687	252	255	260	rVB5	16871	25939	0.30%	0.027%
11	4.987	299	306	316	rVB2	192599	405868	4.73%	0.419%
12	5.081	318	322	329	rVB3	29471	47467	0.55%	0.049%
13	5.869	452	456	459	rBV5	16797	21282	0.25%	0.022%
14	6.281	521	526	532	rVB	44247	65321	0.76%	0.067%
15	6.499	558	563	569	rBV10	10701	19847	0.23%	0.020%
16	7.058	651	658	672	rVB	1445013	2306013	26.88%	2.378%
17	7.228	680	687	695	rBV	1127630	1754363	20.45%	1.809%
18	7.422	714	720	728	rVB	1429728	2228193	25.98%	2.298%
19	7.510	730	735	742	rVB6	15277	26456	0.31%	0.027%
20	7.710	765	769	774	rVB5	11570	18645	0.22%	0.019%
21	7.893	793	800	808	rBV	1527504	2414648	28.15%	2.490%
22	7.952	808	810	816	rVB2	16701	27333	0.32%	0.028%
23	8.610	914	922	931	rBV	1589847	2606722	30.39%	2.689%
24	8.728	936	942	950	rVB	224657	370497	4.32%	0.382%
25	9.063	990	999	1010	rBV	1291047	2150777	25.07%	2.218%
26	9.169	1014	1017	1021	rVV6	13319	20187	0.24%	0.021%
27	9.222	1021	1026	1031	rVB9	12060	21581	0.25%	0.022%
28	9.457	1061	1066	1073	rVB2	34169	58371	0.68%	0.060%
29	9.534	1073	1079	1081	rBV7	13677	25555	0.30%	0.026%
30	9.787	1113	1122	1131	rBV	1095362	1780605	20.76%	1.836%
31	9.946	1144	1149	1155	rBV4	25459	55462	0.65%	0.057%
32	10.004	1155	1159	1165	rVB9	11206	22648	0.26%	0.023%
33	10.152	1178	1184	1189	rBV2	30247	56258	0.66%	0.058%
34	10.328	1207	1214	1224	rBV	1634671	2704878	31.53%	2.790%
35	10.704	1270	1278	1289	rBV	1968668	3333248	38.86%	3.438%
36	10.851	1296	1303	1312	rBV	544787	1004996	11.72%	1.037%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
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37	11.240	1363	1369	1378	rBV5	69089	131316	1.53%	0.135%
38	11.498	1408	1413	1424	rBV4	20794	51014	0.59%	0.053%
39	12.116	1515	1518	1523	rVB7	14607	22958	0.27%	0.024%
40	12.204	1526	1533	1538	rBV5	11871	19152	0.22%	0.020%
41	12.293	1543	1548	1553	rVB2	27246	48039	0.56%	0.050%
42	13.063	1674	1679	1683	rBV7	10558	17193	0.20%	0.018%
43	13.357	1725	1729	1733	rVB4	20452	31157	0.36%	0.032%
44	13.510	1752	1755	1761	rVB7	9343	17201	0.20%	0.018%
45	13.863	1810	1815	1818	rBV7	15144	23757	0.28%	0.025%
46	13.951	1821	1830	1839	rBV	2673680	3612188	42.11%	3.726%
47	14.057	1844	1848	1856	rVB	11470	23837	0.28%	0.025%
48	14.240	1871	1879	1890	rBV	2991265	4446365	51.84%	4.586%
49	14.545	1922	1931	1942	rBV	2783598	4090082	47.68%	4.218%
50	14.751	1960	1966	1977	rBV	604371	1112482	12.97%	1.147%
51	14.963	1997	2002	2008	rVB5	52344	91557	1.07%	0.094%
52	15.269	2051	2054	2059	rVV6	20745	29074	0.34%	0.030%
53	15.316	2059	2062	2068	rVV6	13488	23695	0.28%	0.024%
54	15.381	2071	2073	2079	rVB6	12871	15662	0.18%	0.016%
55	15.539	2093	2100	2107	rVV	3867617	5441758	63.44%	5.613%
56	15.663	2115	2121	2128	rBV	929085	1298154	15.13%	1.339%
57	15.798	2137	2144	2146	rBV	244709	363163	4.23%	0.375%
58	15.822	2146	2148	2155	rVB	240489	288799	3.37%	0.298%
59	15.904	2157	2162	2174	rVB	406247	549887	6.41%	0.567%
60	16.134	2196	2201	2206	rBV	148652	197894	2.31%	0.204%
61	16.216	2211	2215	2217	rBV5	16166	23054	0.27%	0.024%
62	16.251	2217	2221	2225	rVB6	18676	26036	0.30%	0.027%
63	16.357	2236	2239	2245	rBV5	25408	33855	0.39%	0.035%
64	16.410	2245	2248	2251	rBV5	20144	25251	0.29%	0.026%
65	16.469	2255	2258	2262	rVB2	32507	40312	0.47%	0.042%
66	16.528	2263	2268	2273	rBV	60647	89399	1.04%	0.092%
67	16.686	2293	2295	2301	rVB7	14841	18281	0.21%	0.019%
68	16.828	2313	2319	2324	rBV	388724	508771	5.93%	0.525%
69	17.051	2354	2357	2362	rBV7	13222	20987	0.24%	0.022%
70	17.181	2374	2379	2380	rBV3	40119	55088	0.64%	0.057%
71	17.210	2380	2384	2392	rVB	424934	588233	6.86%	0.607%
72	17.298	2392	2399	2408	rBV	3038270	4685467	54.62%	4.833%
73	17.404	2411	2417	2425	rVV	3729612	5329707	62.13%	5.497%
74	17.481	2425	2430	2439	rVB	481138	726865	8.47%	0.750%
75	17.586	2444	2448	2451	rBV2	26827	37616	0.44%	0.039%
76	17.751	2472	2476	2479	rBV3	52223	71815	0.84%	0.074%

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77	17.781	2479	2481	2488	rVB7	35842	52270	0.61%	0.054%
78	17.904	2495	2502	2507	rBV	727340	1245690	14.52%	1.285%
79	17.957	2509	2511	2517	rVB7	21472	27883	0.33%	0.029%
80	18.022	2517	2522	2525	rBV5	46138	90051	1.05%	0.093%
81	18.086	2531	2533	2537	rVB5	41943	48664	0.57%	0.050%
82	18.139	2538	2542	2545	rBV	120394	167450	1.95%	0.173%
83	18.175	2545	2548	2558	rVB	287336	401752	4.68%	0.414%
84	18.351	2575	2578	2585	rBV4	52684	92456	1.08%	0.095%
85	18.410	2585	2588	2591	rVV5	62236	84215	0.98%	0.087%
86	18.451	2591	2595	2601	rVB	190486	241832	2.82%	0.249%
87	18.586	2615	2618	2621	rBV	119640	131360	1.53%	0.135%
88	18.733	2639	2643	2645	rBV2	114093	158312	1.85%	0.163%
89	18.798	2651	2654	2658	rVB2	60178	70673	0.82%	0.073%
90	18.951	2678	2680	2685	rBV2	43902	59731	0.70%	0.062%
91	19.092	2702	2704	2709	rVB2	71595	81404	0.95%	0.084%
92	19.145	2710	2713	2715	rBV2	81741	105074	1.22%	0.108%
93	19.404	2754	2757	2761	rBV4	132422	175403	2.04%	0.181%
94	19.639	2791	2797	2804	rBV	4913299	6566289	76.55%	6.772%
95	20.627	2963	2965	2969	rVB	239098	231336	2.70%	0.239%
96	21.086	3039	3043	3052	rBV	948348	1529430	17.83%	1.577%
97	21.286	3073	3077	3082	rVB	7445874	7680161	89.54%	7.921%
98	21.392	3090	3095	3101	rVB	3798942	4954140	57.76%	5.110%
99	23.680	3478	3484	3498	rVB2	3909161	8577704	100.00%	8.847%
100	23.839	3506	3511	3520	rVB2	2830846	5666671	66.06%	5.845%

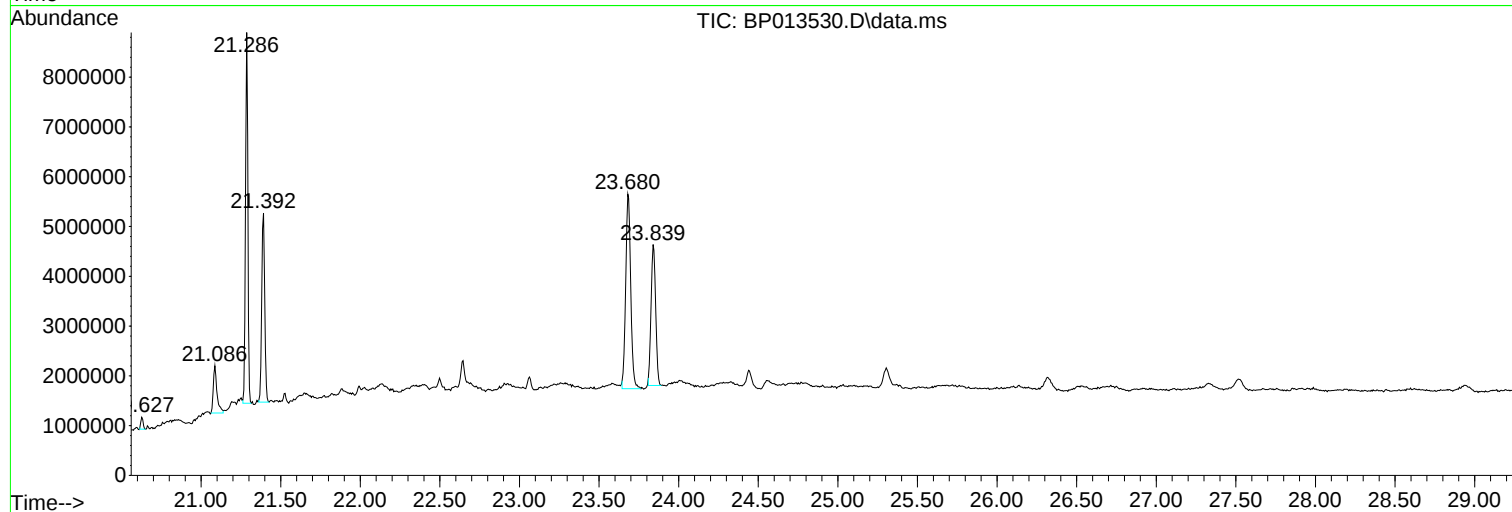
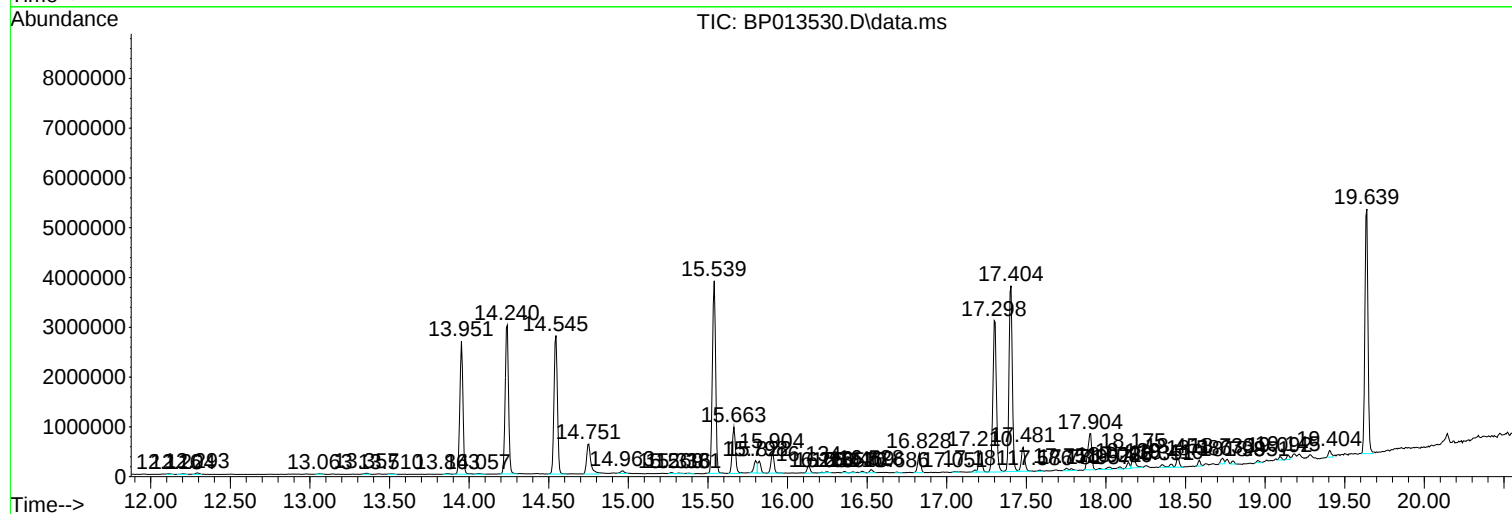
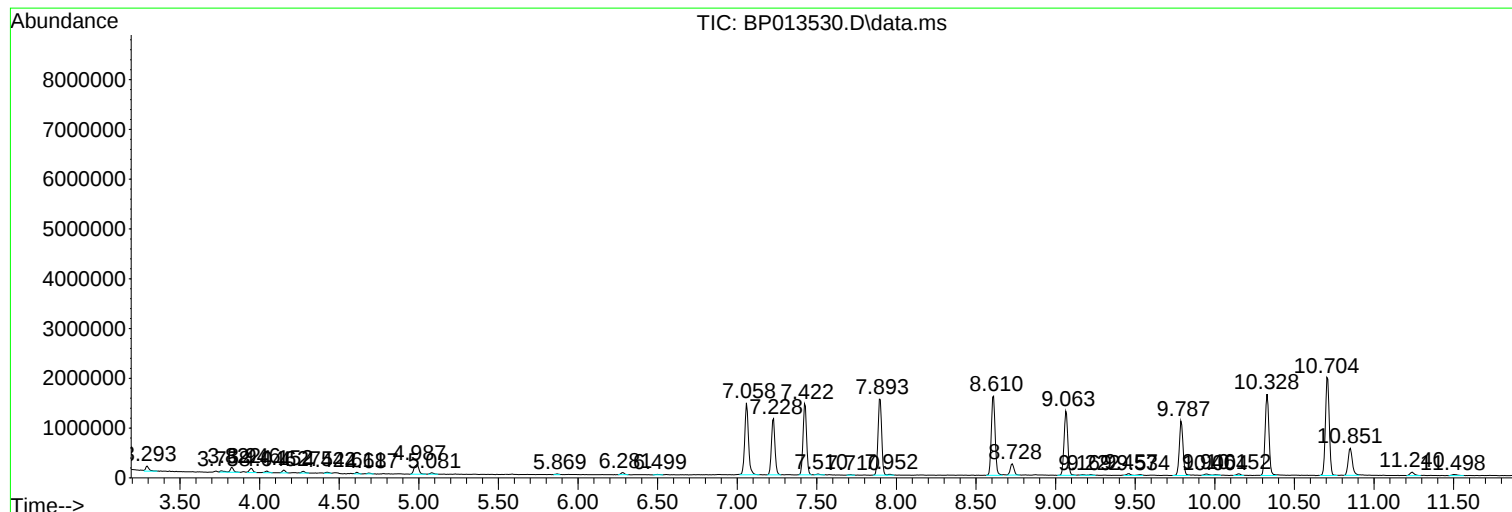
Sum of corrected areas: 96956725

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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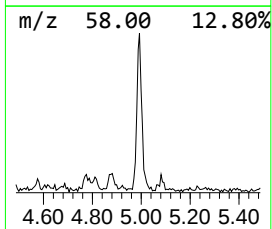
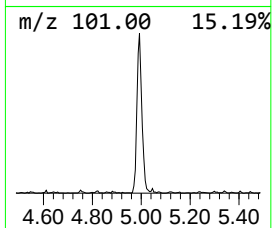
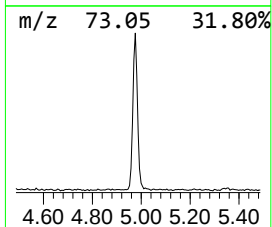
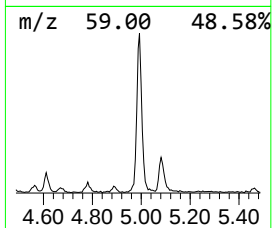
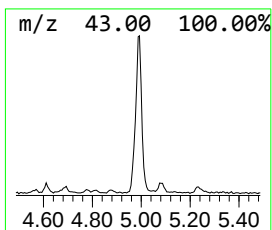
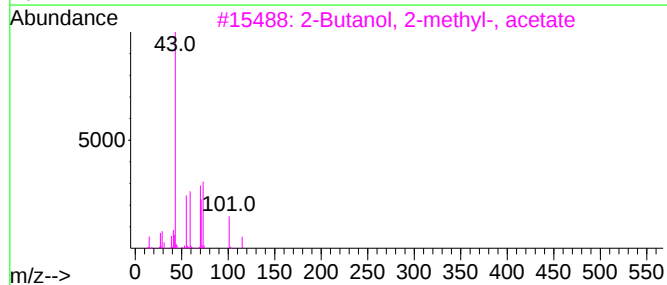
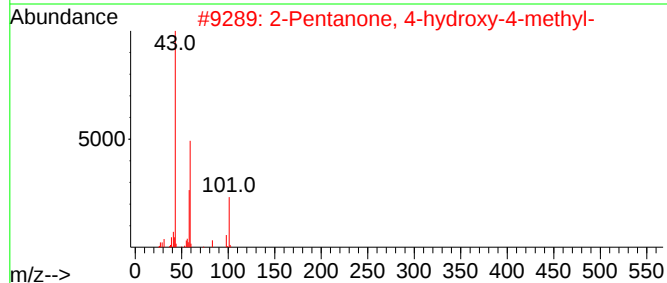
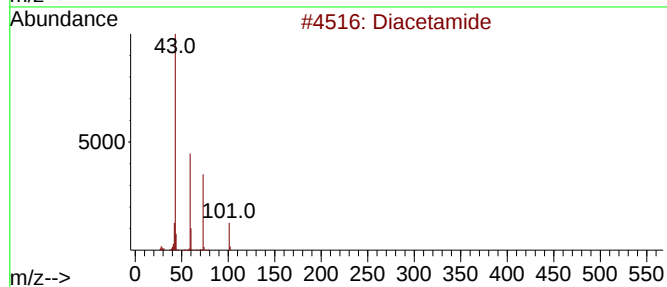
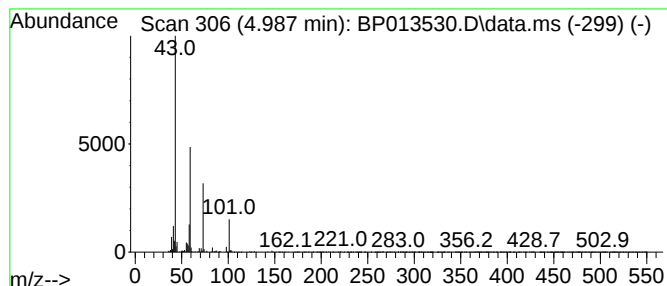
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Diacetamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.987	3.36 ng/ul	405868	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diacetamide	101	C4H7NO2	000625-77-4	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3			2-Butanol, 2-methyl-, acetate	130	C7H14O2	000625-16-1	40
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	33



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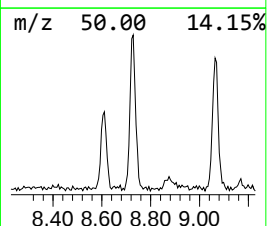
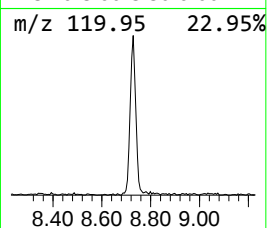
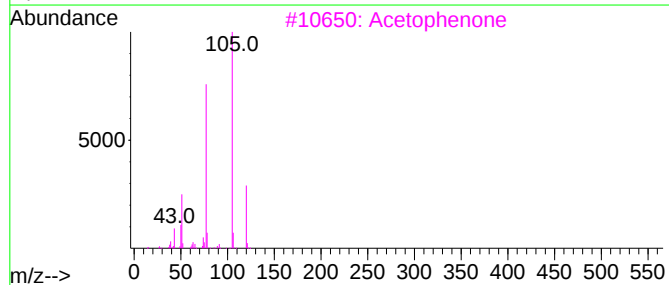
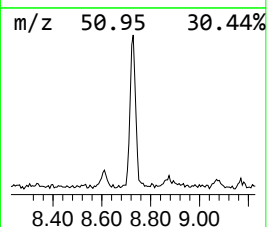
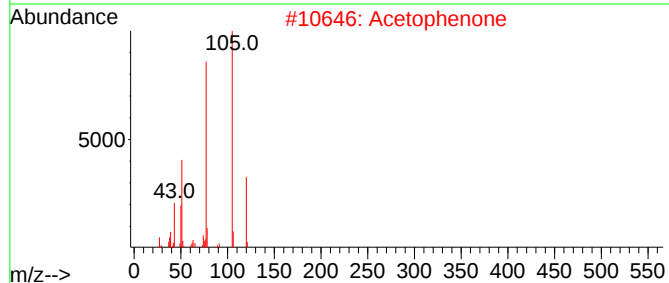
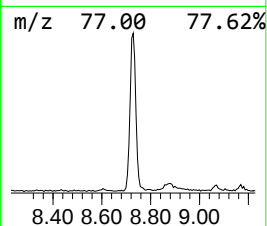
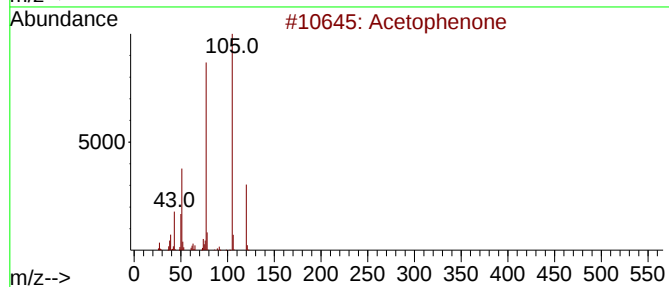
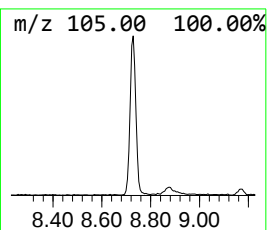
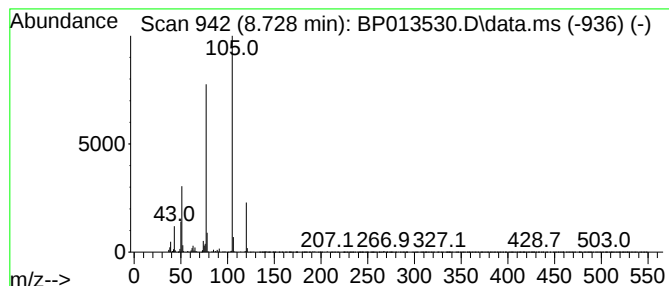
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Acetophe none Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.728	3.07 ng/ul	370497	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	97
2	Acetophenone			120	C8H8O	000098-86-2	94
3	Acetophenone			120	C8H8O	000098-86-2	93
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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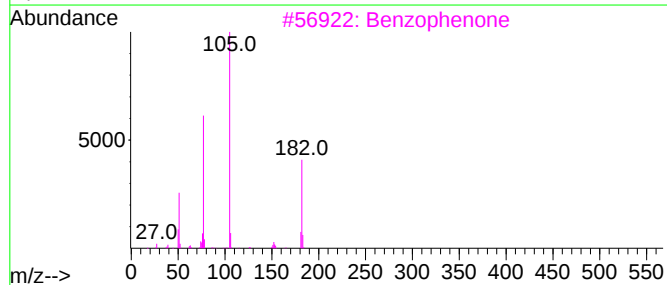
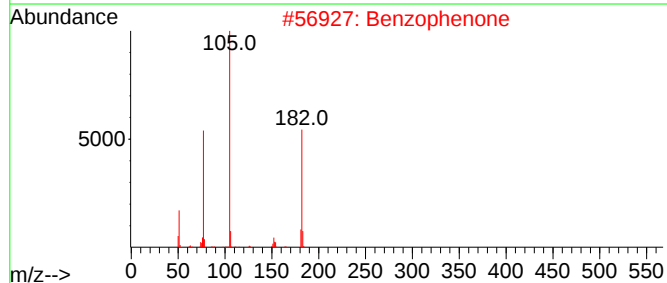
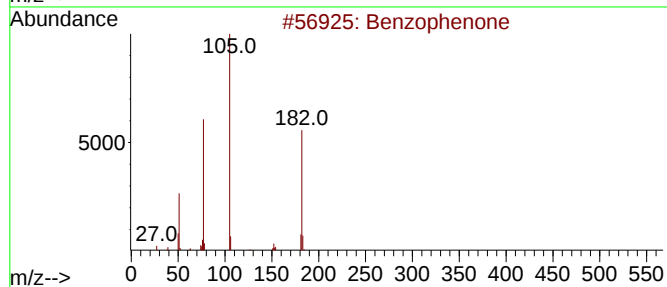
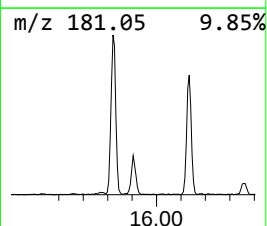
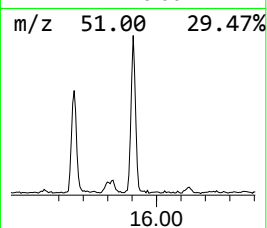
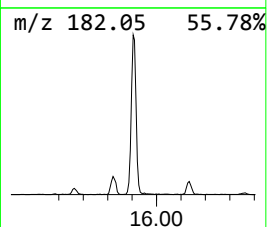
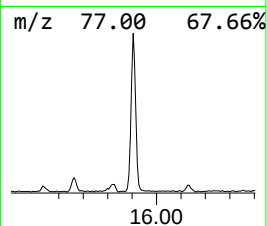
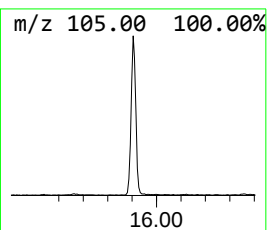
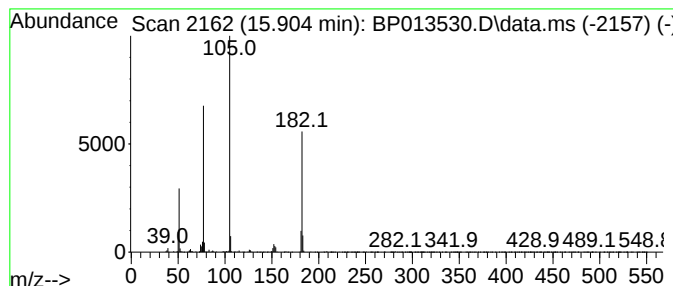
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.904	2.69 ng/ul	549887	Acenaphthene-d10	14.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Z9

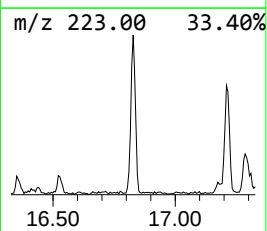
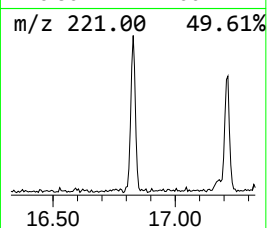
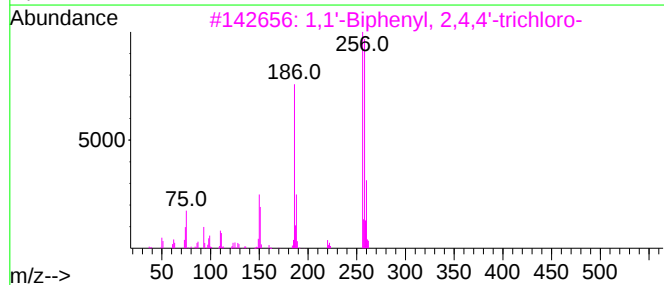
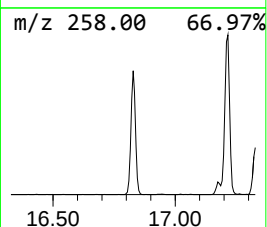
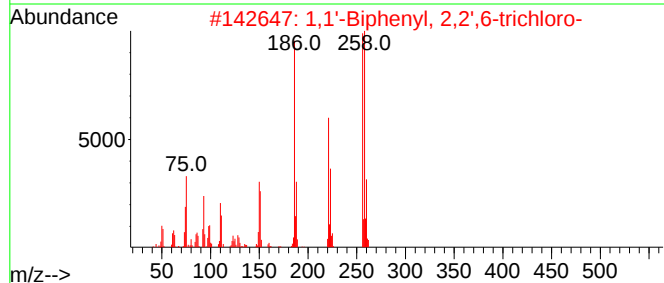
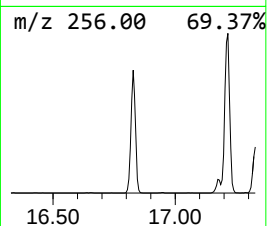
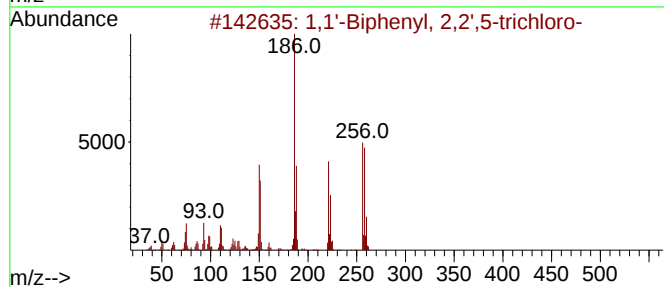
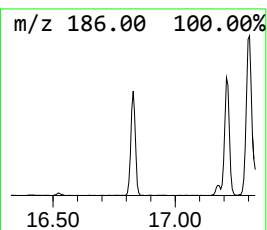
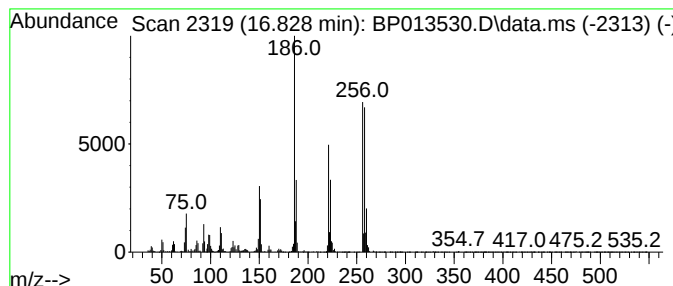
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.828	2.17 ng/ul	508771	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
2	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99		
3	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
4	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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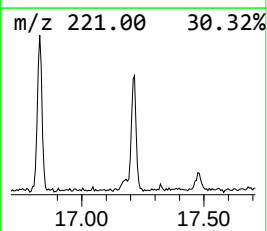
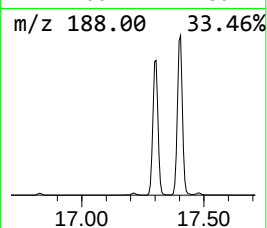
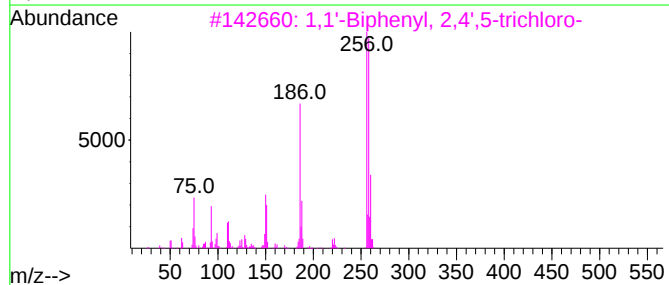
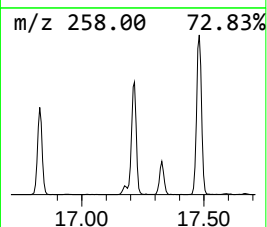
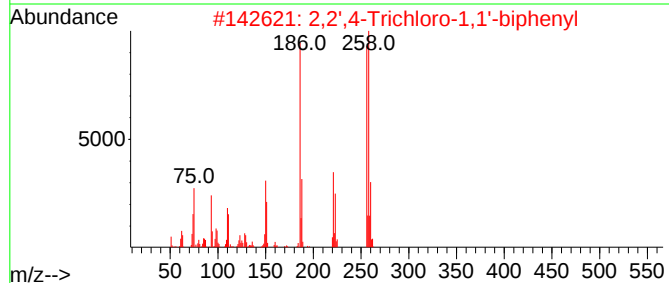
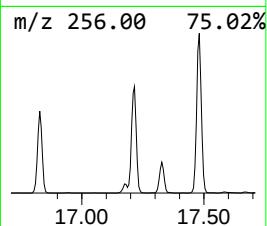
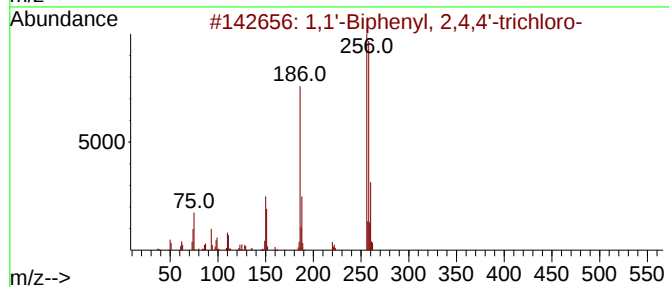
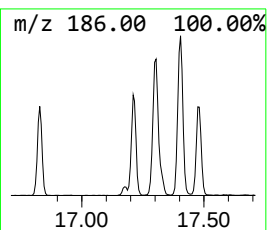
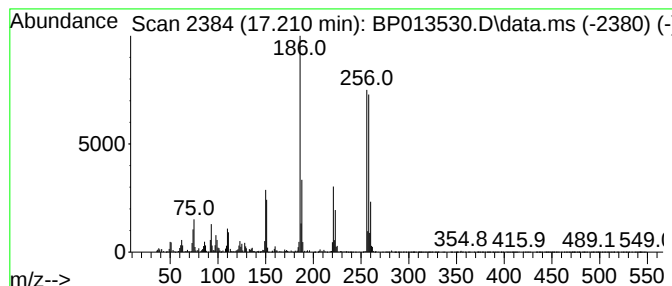
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.210	2.51 ng/ul	588233	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	2,2',4-Trichloro-1,1'-biphenyl	256	C12H7Cl3	037680-66-3	99		
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99		
4	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98		
5	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Z9

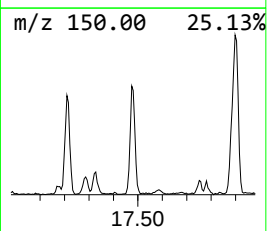
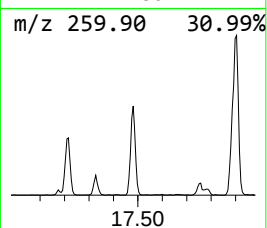
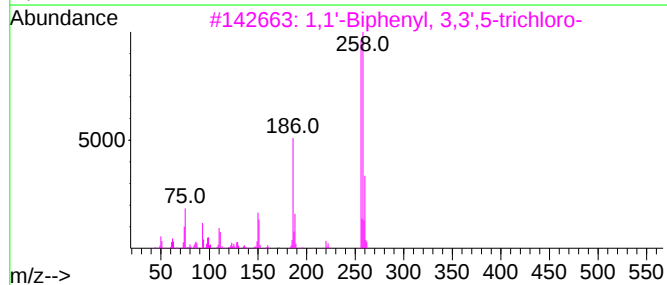
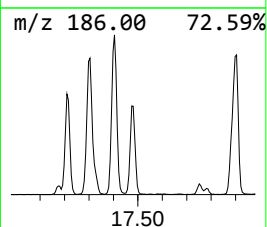
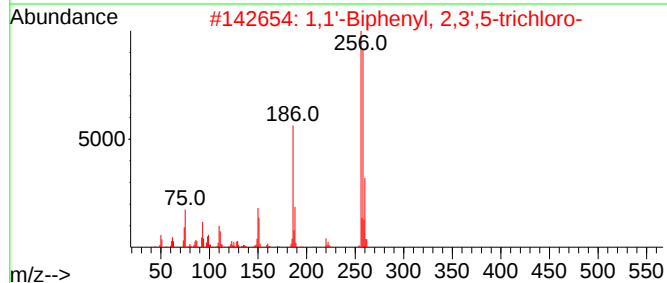
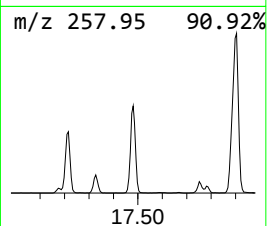
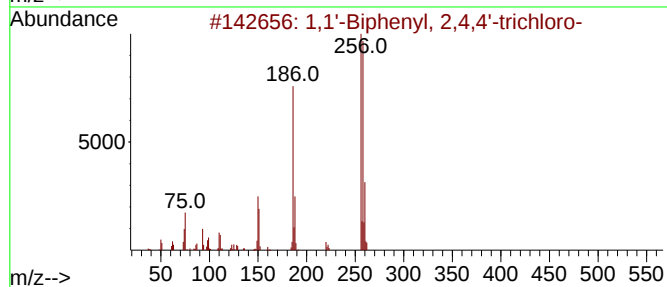
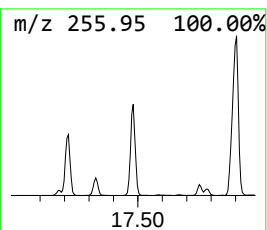
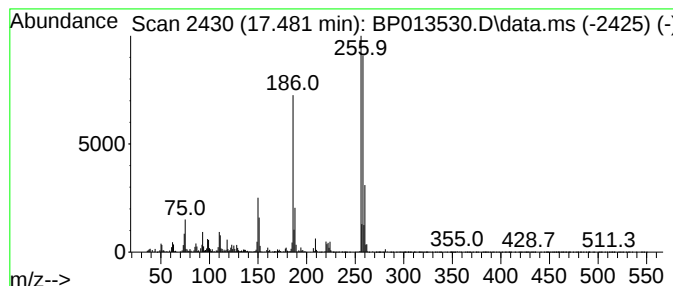
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,1'-Biphenyl, 2,3',5-trich... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.480	3.10 ng/ul	726865	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		
3	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
5	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Z9

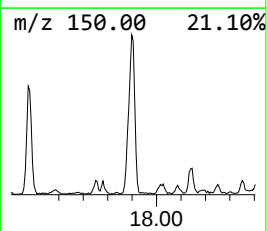
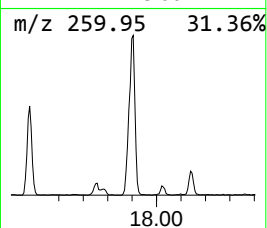
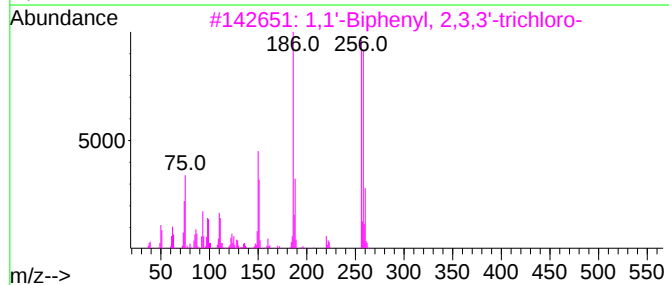
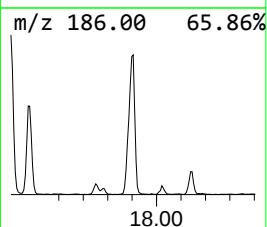
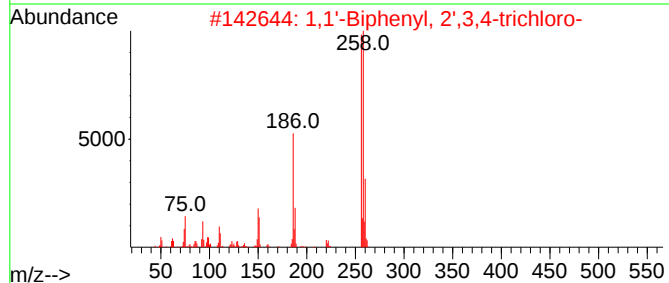
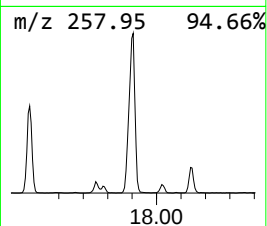
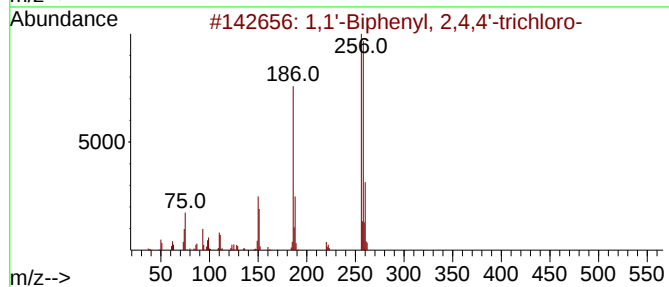
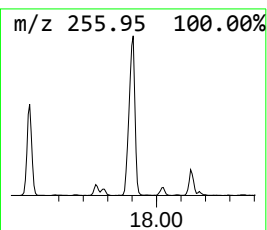
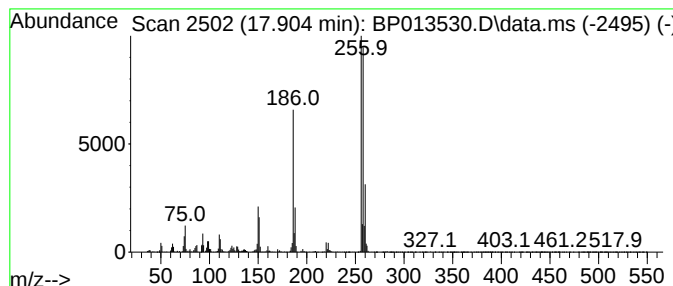
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.904	5.32 ng/ul	1245690	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		
3	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
4	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99		
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

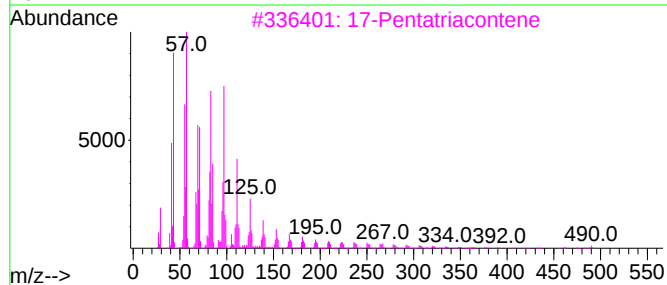
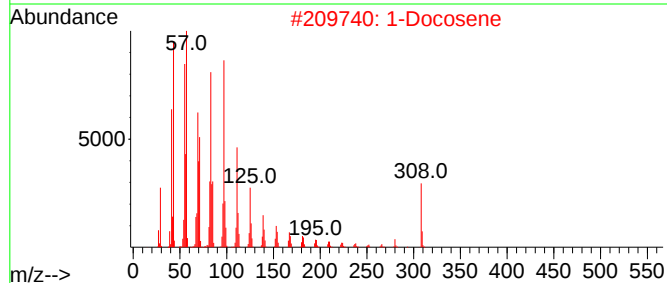
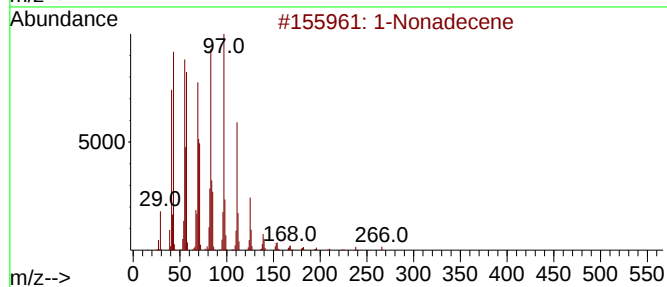
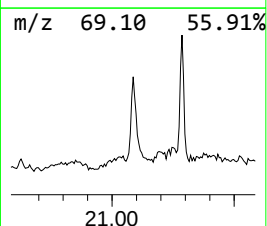
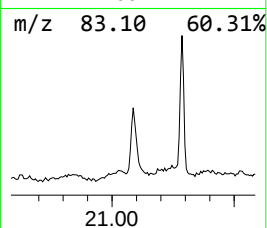
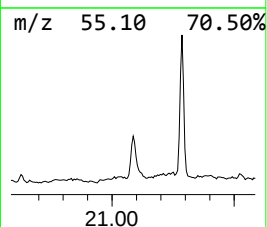
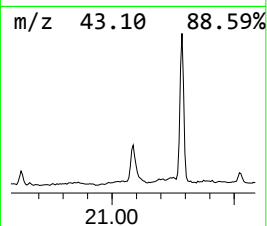
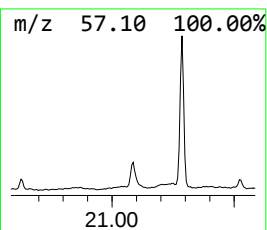
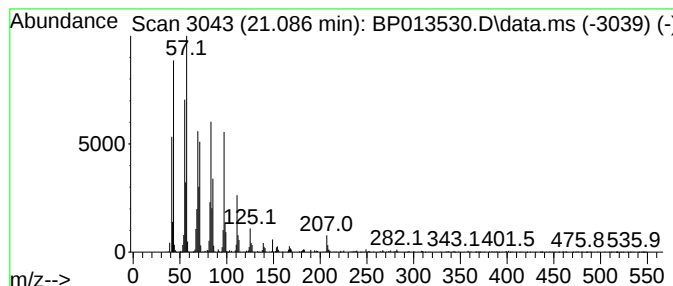
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.086	6.17 ng/ul	1529430	Chrysene-d12		21.392	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Nonadecene		266	C19H38	018435-45-5	96
2	1-Docosene		308	C22H44	001599-67-3	95
3	17-Pentatriacontene		491	C35H70	006971-40-0	95
4	Pentadecafluorooctanoic acid, oc...		666	C26H37F15O2	1000406-04-8	94
5	1-Heneicosyl formate		340	C22H44O2	077899-03-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Acetophe none	8.728	3.1	ng/ul	370497	1	7.899	2414650	20.0
Benzophenone	15.904	2.7	ng/ul	549887	3	14.545	4090080	20.0
1,1'-Biphenyl, ...	16.828	2.2	ng/ul	508771	4	17.304	4685470	20.0
1,1'-Biphenyl, ...	17.210	2.5	ng/ul	588233	4	17.304	4685470	20.0
1,1'-Biphenyl, ...	17.480	3.1	ng/ul	726865	4	17.304	4685470	20.0
1,1'-Biphenyl, ...	17.904	5.3	ng/ul	1245690	4	17.304	4685470	20.0
(DEL) Alkane: S...	21.086	6.2	ng/ul	1529430	5	21.392	4954140	20.0